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**Research Article** 

# GC-MS ANALYSIS OF BIOACTIVE COMPONENTS IN GAULTHERIA FRAGRATISSIMA WALL.

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### ABSTRACT

Medicinal plant species are being used as folk medicine by various tribal and local communities in Tamilnadu as well as all over India since ancient days. The present study of phytochemical analysis of aerial parts of *Gaultheria fragratissima* Wall. was extracted with methanol and analysed using GC-MS. Nearly fifteen bioactive chemical compounds were identified in crude extracts. Predominantly, m-Ethylbenzonitrile, 1-Mehtoxy-1-buten-3-yne, 4-Fluorobenzyl alcohol, 5-Methylcyclopent-1-ene-1-carboxyl, Methyl Salicylate, 2-Hexenal, 2ethyl- and Cyclohexane, 1, 3-dimethyl-2-methyl compounds were present. This analysis revealed that the existence of m-Ethylbenzonitrile and 1-Mehtoxy-1-buten-3-yne as major constituents. These different active phytochemicals have been found to possess a wide range of biological applications.

Keywords: Medicinal plant, GC-MS, Gaultheria fragratissima, Phytochemicals

#### INTRODUCTION

Medicinal plants used for traditional medicine contain a wide range of substances that can be used to treat chronic as well as infectious diseases. Clinical microbiologists have great interest in screening of medicinal plants for new therapeutics (Periyasamy et al., 2010). The therapeutic efficacy of many indigenous plants for various diseases has been described by traditional herbal medicinal practitioners. Even the World Health Organization (WHO) supports the use of medicinal plants, provided it is proven to be efficacious, safe, less toxic, available and reliable natural resource (Lalitha Easwari et al., 2013). The secondary metabolites are alkaloids, phenolics, flavonoids, essential oils and other organic constituents. These secondary metabolites are usually produced in different parts of the plants like roots, leaves, fruits and seeds and then translocated to other parts of plant for storage. Knowledge about these medicinally active constituents makes their application as medicine as mentioned in the various pharmacopeias.

The genus Gaultheria comprises about 200 species which are native to wide geographical areas ranging from Andes, North America, Australia and nearby islands to South-East Asia. In India, it is found mainly in the eastern Himalayas, it is found throughout the North-Eastern region including the Indo-Nepal and Indo-Bhutan territories at an altitude between 1,800 and 2,500 m whereas in Western Ghats, it is found in Nilgiri, Palani Hills of Tamilnadu up to an altitude of 1,500 m with low frequency with low frequency and density (Bantawa et al., 2011). The plant has been used as an antiseptic, carminative, flavouring agent (confectionaries, herbal tea. toothpaste, etc.) and condiment and also in rheumatic and arthritis treatments. The oil has a powerful antiseptic property and is a constituent for several insecticidal and insect repellent preparations (Ranyaphi et al., 2012). Hence during the present investigations phytochemical screening of Gaultheria fragrantissima is carried on with a view to analyse the presence of chemical constituents that secondary metabolites, with a view to recommend their application in pharmaceutical industry.

#### Materials and methods

### **Plant materials**

The plant samples were collected from Coimbatore District, Tamilnadu. The botanical identify of the plant was confirmed by Dr. S. Padmavathy, Professor, PG & Research Department of Botany, Nirmala College for Women, Coimbatore, Tamilnadu.

#### Plant sample preparation

50gm powdered plant material was soaked in 200 ml of methanol overnight and then centrifuged through Whatman no.1 filter paper along with 2gm sodium sulfate to remove the sediments and traces of water in the filtrate. Before filtering the filter paper along with sodium sulphate is wetted with methanol. The filtrate is then concentrated by bubbling nitrogen gas into the solution and reduces the volume to 1ml. The extract contains both polar and non-polar phytocomponents.

# **GC-MS** analysis

GC-MS analysis was carried out on a GC clarus 500 Perkin Elmer system comprising a AOC-20I auto sampler and gas chromatograph interfaced to a mass spectrophotometer instrument employing the following conditions: Column Elite-1 fused silica capillary column (30mm×0.25mm I.D ×1  $\mu$  M df, composed of 100% Dimethyl poly siloxane ), operating in electron impact mode at 70 eV; helium (99.999%) was used as carrier gas at a constant flow of 1ml/min and an injection volume of 0.5  $\mu$  l was employed (split ratio of10:1) injector temperature 250 °C; ion-source temperature 280 °C. The oven temperature was programmed from 110 °C (isothermal for 2 min), with an increase of 10 °C/min, to 200 °C, then 5 °C/min to 280 °C, ending with a 9 min isothermal at 280 °C. Mass spectra were taken at 70 eV; a scan interval of 0.5 seconds and fragments from 45 to 450 Da. Total GC running time is 46min.

#### **Results and Discussion**

Medicinal plants are being used as valuable sources of food and medicine for the prevention of illness and maintenance of human health. In India, many indigenous plants are widely consumed as food or home remedies especially in the treatment or management of common diseases. The composition and identification of the main components present in the Gaultheria fragrantissima are shown in (Table 1). Fifteen compounds were identified in G. fragrantissima by GC-MS analysis. The active principles with their retention time (RT), percentage area and Reference No. are presented in (Table 1 and Fig 1). It was found that the main constituents of leaves m-Ethylbenzonitrile (22.57%), 1-Mehtoxy-1-buten-3-yne (21.49%), 4-Fluorobenzyl alcohol (9.83%), 5-Methylcyclopent-1-ene-1-carboxyl (8.23%), Methyl Salicylate (7.17%), 2-Hexenal 2ethyl- (6.43%), 1,3-dimethyl-2-methyl (6.05%),2,4-Decadienal Cyclohexane. (4.80%), N-(n-Propyl)acetamide (3.41%).

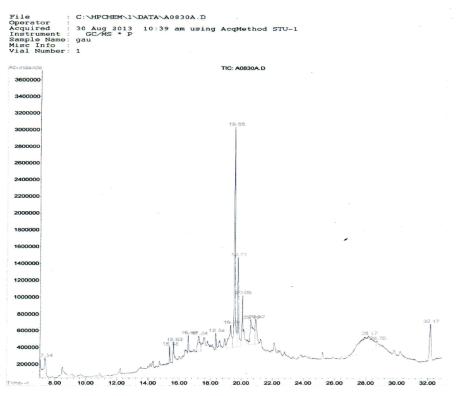


Figure 1: GC-MS studies of Gaultheria fragrantissima Wall.

Gaultherin is a monotropitin a conjugated compound made up of methyl salicylate and disaccharides of glucose and xylose, called primevorose. Methyl salicylate is the natural precursor of pharmaceutical aspirin. In the US, there are more than 40 products that contain methyl acetate as an active ingredient to treat various kinds of external pains. The oil has high demand in pharmaceutical and perfumery industries (Ranyaphi *et al.*, 2012). In the present study fifteen chemical constituents have been identified from methanol extract of aerial parts of *Gaultheria fragrantissima* by GC-MS analysis. The existence of various bioactive chemical compounds proves the use of this plant various ailment by traditional medical practitioners.

Table 1: GC-MS studies of Gaultheria	fragrantissima Wall.
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Pk#	RT	Area%	Library/ID	Ref#	CAS# Qu	al
1	7.34	1.98	C: DATABASE NIST98.L		and the set of the set of	
			Isopropyl Alcohol		000067-63-0	
			Isopropyl Alcohol	110736	000067-63-0	64
			Propylene Glycol	11460	000057-55-6	64
2	15.37	1.20	C: DATABASE NIST98 L			
			Urea, trimethyl-		000632-14-4	
			Tetraacetyl-d-xylonic nitrile		1000130-04-4	
			2-Formylhistamine	10720	1000132-95-8	40
3	15.61	2 25	C:\DATABASE\NIST98.L			
5 15.	10.01		4H-Pyran-4-one, 2.3-dihydro-3.5-di	109420	028564-83-2	87
			4H-Pyran-4-one, 2.3-dihydro-3.5-di	4836	028564-83-2	62
			2-Tetrazene. 1.1-diethyl-4.4-dimet	11216	014866-81-0	43
4	16.56	1.48	C:\DATABASE\NIST98.L 2-Furancarboxaldehyde. 5-(hydroxym	118306	000067-47-0	60
			2-Furancarboxaldehyde, 5-(hydroxym	118397	000067-47-0	53
			4-Mercaptophenol	57293	000637-89-8	35
5	17.24	3.41	C:\DATABASE\NIST98.L 4H-Pyran-4-one, 2,3-dihydro-3,5-di	4836	028564-83-2	64
			4H-Pyran-4-one, 2,3-dihydro-3,5-di	109420	020564 83 2	59
			N-(n-Propyl)acetamide	10694	005331-48-6	27
			-			
6	18.34	1.27	C:\DATABASE\NIST98.L N-Methoxy-1-ribofuranosyl-4-carbox	10000	1000129 57 7	5
			N-Methoxy-1-riboluranosy1-4-carbox	10000	1000130-04-4	1 4
			Tetraacetyl-d-xylonic nitrile		1000225-93-3	
			N-(Glycyl)alanine	10524	1000220-00	
7	19.29	6.43	C: DATABASE NIST98.L			
					000645-62-5	
			Thiophene-3-boronic acid		006165-69-1	
			2-Hexenal, 2-ethyl-	111278	000645-62-5	20
8	19.55	21.49	C:\DATABASE\NIST98.L			120125
			2-Furancarboxaldehyde. 5-(hydroxym	118397	000067-47-0	90
			Thiocyanic acid, 2-propynyl ester		024309-48-6	
			1-Methoxy-1-buten-3-yne	1343	002798-73-4	47
9	19.76	9.83	C:\DATABASE\NIST98.L			
9 1.			2-Eurancarboyaldebyde 5-(bydroxym	118397	000067-47-0	91
			2-Furancarboxaldehyde. 5-(hydroxym	118306	000067-47-0	55
			4-Fluorobenzyl alcohol	42355	000459-56-3	50
10	20 06	7 17	C: DATABASE NIST98.L			
10	20.00	1.71	Methyl Salicylate		000119-36-8	
			Methyl Salicylate	120910	000119-36-8	96
			Methyl Salicylate		000119-36-8	
	C					
11	20.60	8.23	C:\DATABASE\NIST98.L 2-Furancarboxaldehyde, 5-(hydroxym	118307	000067-47-0	46
	2004		2-Furancarboxaldehyde, 5-(hydroxym 2-Furancarboxaldehyde, 5-(hydroxym	118306	000067-47-0	46
			5-Methylcyclopent-1-ene-1-carboxyl	32139	053623-50-0	3.8
			2-MeenArcAcrobene-r-ene-r-carpoxAr			

```
12
          4.80 C:\DATABASE\NIST98.L
                    2.4-Decadienal
2.4-Decadienal, (E.E)-
                                                                               116043 002363-88-4 38
116039 025152-84-5 38
                    2,4-Decadienal, (E.E)-
                                                                                 31387 025152-84-5 35
28.16 22.57 C:\DATABASE\NIST98.L
4-Cyclopropylmethylbenzonitrile
6-Methyl benzobicyclo[2.2.1] hepte
m-Ethylbenzonitrile
                                                                                52135 1000110-39-3 35
52134 140224-88-0 32
                                                                                 52140 1000118-64-1 30
28.70 1.85 C:\DATABASE\NIST98.L
6-Methyl benzobicyclo[2.2.1] hepte
4-Cyclopropylmethylbenzonitrile
                                                                                 52134 140224-88-0 38
                                                                                 52135
                                                                                          1000110-39-3 35
                    Heptadecanoic acid, 3-oxo-, methyl
                                                                                 52116 054934-65-5 27
32.17 6.05 C:\DATABASE\NIST98.L
                   Bicyclo[3.1.1]heptane, 2.6.6-trime 2052 004795-86-2 87
Bicyclo[3.1.1]heptane, 2.6.6-trime 13368 006876-13-7 86
Cyclohexane, 1.3-dimethyl-2-methyl 113932 019781-47-6 43
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